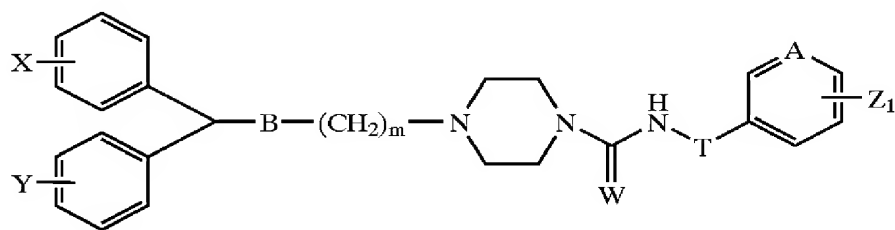


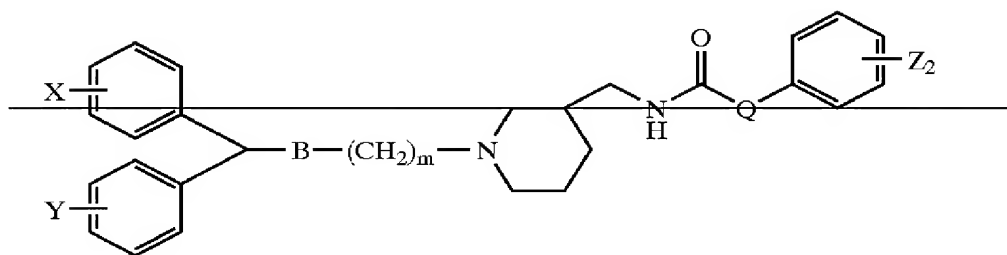
IN THE CLAIMS

1. A compound of Formula I having high affinity for a dopamine transporter ~~having a formula selected from the group consisting of:~~

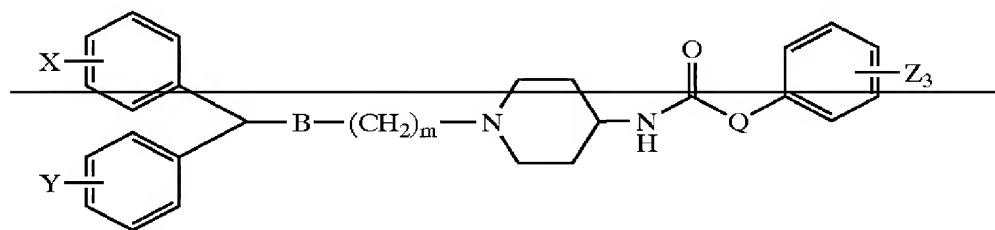
Formula I



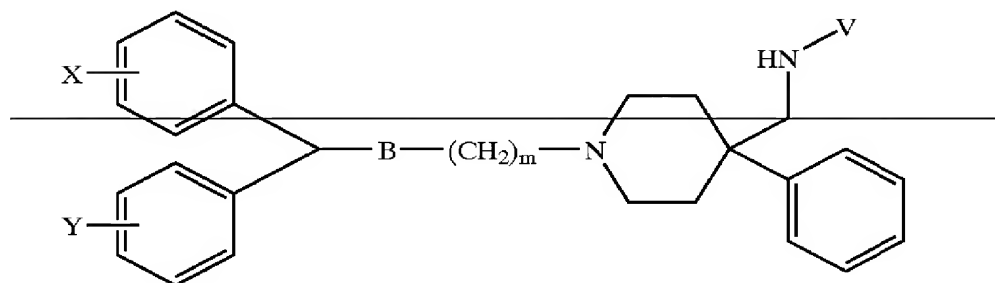
Formula II



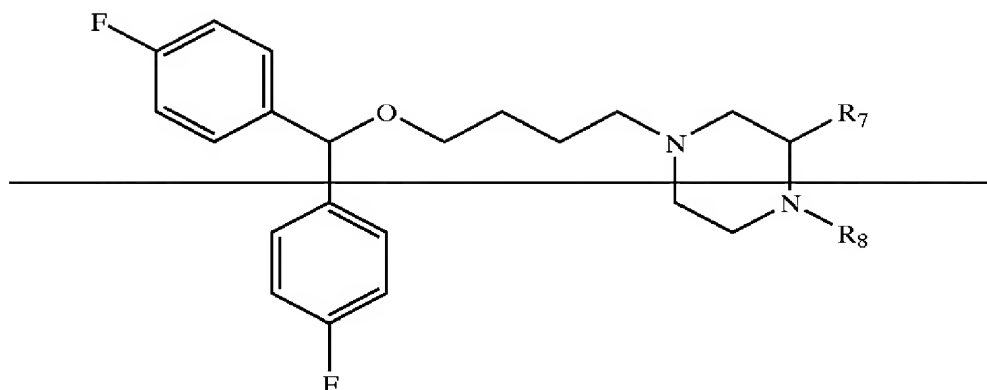
Formula III



Formula IV



Formula V



wherein:

B is -O-, -N(H)-, -C(=O)N(H)-, or -N(H)C(=O)-;

[[n]] m is an integer of 1 to 6; ~~X, Y, Z<sub>1</sub> and Z<sub>2</sub> can be the same or different and are hydrogen, halo, haloalkyl, alkyl, aryl, (C<sub>4</sub>-C<sub>6</sub>) alkoxy, N-alkyl, (C<sub>2</sub>-C<sub>6</sub>) acyloxy, N-alkylene, -SH, -SR, wherein R is from the same group as R<sub>1</sub> and R<sub>2</sub> and can be the same or different than R<sub>1</sub> and R<sub>2</sub>, amino, nitro, cyano, hydroxy, C(=O)OR<sub>6</sub>, C(=O)NR<sub>5</sub>R<sub>6</sub>, NR<sub>3</sub>R<sub>2</sub> or S(=O)<sub>k</sub>R<sub>1</sub> wherein k is 1 or 2, and R<sub>1</sub> to R<sub>6</sub> are independently hydrogen or (C<sub>4</sub>-C<sub>6</sub>) alkyl;~~

X is hydrogen, halo, haloalkyl, alkyl, aryl, C<sub>1</sub>-C<sub>6</sub> alkoxy, N-alkyl, C<sub>2</sub>-C<sub>6</sub> acyloxy, or N-alkylene;

Y is hydrogen, halo, haloalkyl, alkyl, aryl, C<sub>1</sub>-C<sub>6</sub> alkoxy, N-alkyl, C<sub>2</sub>-C<sub>6</sub> acyloxy, or N-alkylene;

Z<sub>1</sub> is hydrogen, halo, haloalkyl, alkyl, aryl, C<sub>1</sub>-C<sub>6</sub> alkoxy, N-alkyl, C<sub>2</sub>-C<sub>6</sub> acyloxy, or N-alkylene;

~~R<sub>1</sub> and R<sub>2</sub> can be the same or different and are hydrogen, (C<sub>4</sub>-C<sub>6</sub>) alkyl, hydroxyalkyl or mercaptoalkyl, C(=O)OR<sub>1</sub>, cyano, (C<sub>4</sub>-C<sub>6</sub>) alkenyl, (C<sub>2</sub>-C<sub>6</sub>) alkynyl, or 1,2,4-oxadiazol-5-yl optionally substituted at the 3-position by Z<sub>2</sub>, wherein any (C<sub>4</sub>-C<sub>6</sub>) alkyl, (C<sub>4</sub>-C<sub>6</sub>) alkanoyl, (C<sub>2</sub>-C<sub>6</sub>) alkenyl or (C<sub>2</sub>-C<sub>6</sub>) alkynyl can optionally be substituted by 1, 2 or 3 Z<sub>2</sub>; Z<sub>4</sub> is (C<sub>4</sub>-C<sub>6</sub>) alkyl or phenyl, optionally substituted by 1, 2 or 3 Z<sub>4</sub>~~

~~R<sub>7</sub> can be hydrogen, O or phenyl~~

~~R<sub>8</sub> can be hydrogen, phenyl, halophenyl, nitrophenyl, pyridyl, piperonyl or sulfoxonitrophenyl~~

W is O or S<sub>2</sub>

~~T is amino or C<sub>1</sub>-C<sub>6</sub> aminoalkyl~~

A is N or C<sub>1</sub> and

T is C<sub>1</sub>-C<sub>6</sub> alkyl or sulfonyl and

~~V is alkyl (C<sub>0</sub>-C<sub>6</sub>), alkenyl, alkynyl, haloaryl, alkyl phenol, alkyl halophenyl, and R<sub>1</sub> or R<sub>2</sub> as indicated above and~~

~~φ is phenyl, naphthyl, thienyl or pyridinyl.~~

Claims 2-13 **(canceled)**

14. **(new)** The compound of claim 1, wherein X is halo.
15. **(new)** The compound of claim 1, wherein X is *para*-F.
16. **(new)** The compound of claim 1, wherein Y is halo.
17. **(new)** The compound of claim 1, wherein Y is *para*-F.
18. **(new)** The compound of claim 1, wherein B is O.
19. **(new)** The compound of claim 1, wherein m is 4.
20. **(new)** The compound of claim 1, wherein W is O.
21. **(new)** The compound of claim 1, wherein T is C<sub>1</sub>-C<sub>6</sub> alkyl.
22. **(new)** The compound of claim 1, wherein T is -CH<sub>2</sub>-.
23. **(new)** The compound of claim 1, wherein T is sulfonyl.
24. **(new)** The compound of claim 1, wherein A is C.
25. **(new)** The compound of claim 1, wherein Z<sub>1</sub> is halo.
26. **(new)** The compound of claim 1, wherein Z<sub>1</sub> is *para*-F.

27. (new) The compound of claim 1, wherein the compound is selected from the group

